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IMPROVEMENT OF FINITE ELEMENT MESHES: HEAT TRANSFER IN AN INFINITE CYLINDER

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SUMMARY

1 This paper presents an extension of a structural finite element mesh improvement technique to heat conduction analysis. The mesh improvement concept was originally presented by Prager in studying tapered, axially loaded bars. It was further developed by Kittur, et al., who showed that an improved mesh can be obtained by minimizing the trace of the stiffness matrix. In this paper these procedures are extended and applied to the analysis of heat conduction in an infinitely long hollow circular cylinder.

NOMENCLATURE

N^e	shape function for element e
n	number of elements
q	temperature gradient, $^{\circ}\text{C}/\text{mm}$
r	radius, mm
r_o	inner radius of the cylinder, mm
r_n	outer radius of the cylinder, mm
r_{e-1}	inner radius of element e , mm
r_e	inner radius of element $e+1$ or outer radius of element e , mm
T_o	temperature at the inner radius, $^{\circ}\text{C}$
T_n	temperature at the outer radius, $^{\circ}\text{C}$
$[k_{ij}^e]$	stiffness matrix of element e
κ	thermal conductivity, $\text{W}/\text{mm } ^{\circ}\text{C}$

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κ_e thermal conductivity of the material of element e , W/mm °C
 τ trace of the global stiffness matrix

INTRODUCTION

The finite element analyst is continually confronted with the decision of selecting a "good" mesh for a given problem. Unfortunately there seems to be no simple mesh optimization method which is generally applicable. Recently Kittur, et al., (ref. 1) have proposed a method which provides an improved mesh over a uniform mesh for a broad class of structural problems. The method is based upon the simple procedure of node relocation to minimize the trace of the stiffness matrix. Figure 1 presents a flowchart illustrating the procedure.

An advantage of this method is that the analyst may select the nodal positions prior to solving the equilibrium equations. A posteriori methods such as those of Shephard (ref. 2) and Carroll (ref. 3) can then be used to refine the mesh further.

Kittur, et al., have shown that the minimum trace method produces excellent results for the problem studied by Prager (ref. 4): a uniformly tapered, axially loaded bar. This problem is governed by Poisson's equation with Dirichlet boundary conditions.

In this paper we extend and apply the method to analyze the heat conduction through the cross section of an infinitely long hollow circular cylinder. For this problem, the governing equation is also of the form of Poisson's equation. In this case, however, since no heat sources are considered the governing equation reduces to Laplace's equation. Both Dirichlet and mixed boundary conditions are considered. The results are compared with analytical solutions and with finite element solutions obtained with a uniform mesh.

CONFIGURATION AND PROBLEM DEFINITION

Consider a hollow circular cylinder of infinite length having inner and outer radii: r_o and r_n . Let the thermal conductivity be κ , and the temperatures at the inner and outer radii be T_o and T_n respectively. The governing equation for the temperature distribution along a radial line is:

$$\frac{d}{dr} \left[r\kappa \frac{dT}{dr} \right] = 0 \quad (1)$$

The boundary conditions are:

$$\begin{aligned} T &= T_o \quad \text{at} \quad r = r_o, \\ \text{and} \\ T &= T_n \quad \text{at} \quad r = r_n \end{aligned} \quad (2)$$

The solution of equation (1) subject to equation (2) is:

$$T = T_n + (T_o - T_n) \frac{\ln\left(\frac{r_n}{r}\right)}{\ln\left(\frac{r_n}{r_o}\right)} \quad (3)$$

Next, suppose that the temperature gradient at the inner surface is specified as: q_0 . The boundary conditions are then:

$$\begin{aligned} \frac{dT}{dr} &= q_0 \quad \text{at} \quad r = r_0 \\ \text{and} \\ T &= T_n \quad \text{at} \quad r = r_n \end{aligned} \quad (4)$$

In this case the solution of equation (1) is:

$$T = T_n - r_0 q_0 \ln\left(\frac{r_n}{r}\right) \quad (5)$$

FINITE ELEMENT FORMULATION AND MESH OPTIMIZATION

Figure 2 shows the finite element model. It consists of a series of annular elements. For element (e) let the inner and outer radii be r_e and r_{e+1} . The entries of the element stiffness matrix are:

$$[k_{ij}^e] = 2\pi\kappa_e \int_{r_e}^{r_{e+1}} r \frac{dN_i^e}{dr} \frac{dN_j^e}{dr} dr \quad (6)$$

where κ_e is the element conductivity constant and where the element shape functions N_1^e and N_2^e are:

$$N_1^e = \frac{(r_{e+1} - r)}{(r_{e+1} - r_e)} \quad \text{and} \quad (7)$$

$$N_2^e = \frac{(r - r_e)}{(r_{e+1} - r_e)}$$

By carrying out the indicated operations the element stiffness matrix becomes:

$$[k_{ij}^e] = S_e \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} \quad (8)$$

where S_e is defined as:

$$S_e = \pi\kappa_e \left(\frac{r_{e+1} + r_e}{r_{e+1} - r_e} \right) \quad (9)$$

Hence the trace τ of the global stiffness matrix is:

$$\tau = 2 \sum_{e=1}^n S_e \quad (10)$$

where n is the number of elements.

The trace may be minimized with respect to the nodal coordinates by setting the partial derivative of τ with respect to r_e equal to zero and solving for the ratio r_{e+1}/r_e :

$$\frac{\partial \tau_e}{\partial r_e} = 0 \quad (11)$$

Since the derivative of a sum is the sum of the derivatives, and assuming that the conductivity κ is uniform through the body (i.e.: $\kappa_{e-1} = \kappa_e = \kappa_{e+1}$), (11) becomes:

$$\frac{\partial}{\partial r_e} \left[\dots + \frac{r_e + r_{e-1}}{r_e - r_{e-1}} + \frac{r_{e+1} + r_e}{r_{e+1} - r_e} + \dots \right] = 0 \quad (12)$$

which simplifies to:

$$\frac{-2r_{e-1}}{(r_e - r_{e-1})^2} + \frac{2r_{e+1}}{(r_{e+1} - r_e)^2} = 0 \quad (13)$$

Inverting (13) and factoring:

$$\frac{r_e^2 \left(\frac{r_{e+1}}{r_e} - 1 \right)^2}{r_{e+1}} = \frac{r_{e-1}^2 \left(\frac{r_e}{r_{e-1}} - 1 \right)^2}{r_{e-1}} \quad (14)$$

Rearranging:

$$\left(\frac{r_e}{r_{e-1}} \right) \left(\frac{r_{e+1}}{r_e} - 1 \right)^2 = \left(\frac{r_{e+1}}{r_e} \right) \left(\frac{r_e}{r_{e-1}} - 1 \right)^2 \quad (15)$$

Let $A = \frac{r_{e+1}}{r_e}$ and $B = \frac{r_e}{r_{e-1}}$, then (15) becomes:

$$B(A - 1)^2 = A(B - 1)^2 \quad (16)$$

or:

$$AB(A - B) - (A - B) = 0 \quad (17)$$

thus:

$$(A - B)(AB - 1) = 0 \quad (18)$$

Equation (18) implies that either $(A - B) = 0$ or $(AB - 1) = 0$. If $(AB - 1) = 0$, then $AB = 1$, which means $\frac{r_{e+1}}{r_e} \frac{r_e}{r_{e-1}} = 1 = \frac{r_{e+1}}{r_{e-1}}$ which is impossible. Therefore $(A - B) = 0$, which implies that $A = B$. This leads to the relatively simple relation:

$$\frac{r_{e+1}}{r_e} = \frac{r_e}{r_{e-1}} = \gamma \quad (19)$$

where γ is a constant. The optimal nodal positions are found through repeated use of equation (19):

$$\frac{r_2}{r_1} = \frac{r_1}{r_0} = \frac{r_3}{r_2} = \frac{r_2}{r_1} = \frac{r_{e+1}}{r_e} = \frac{r_e}{r_{e-1}} = \frac{r_n}{r_{n-1}} = \frac{r_{n-1}}{r_{n-2}} = \gamma \quad (20)$$

Then:

$$\frac{r_n}{r_0} = \left(\frac{r_n}{r_{n-1}}\right)\left(\frac{r_{n-1}}{r_{n-2}}\right) \cdot \cdot \cdot \left(\frac{r_{e+1}}{r_e}\right)\left(\frac{r_e}{r_{e-1}}\right) \cdot \cdot \cdot \left(\frac{r_3}{r_2}\right)\left(\frac{r_2}{r_1}\right)\left(\frac{r_1}{r_0}\right) \quad (21)$$

$$\frac{r_n}{r_0} = \gamma \cdot \gamma \cdot \cdot \cdot \gamma \cdot \gamma \cdot \cdot \cdot \gamma \cdot \gamma \cdot \gamma = \gamma^n \quad (22)$$

thus: $r_n = r_0 \gamma^n$ and $\gamma = (r_n/r_0)^{1/n}$. Changing n to e , we have:

$r_e = r_0 \gamma^e$ or:

$$r_e = r_0 \left(\frac{r_n}{r_0}\right)^{e/n} \quad (23)$$

NUMERICAL EXAMPLES

Example 1

To illustrate the effectiveness of the method consider an annular cylinder with the following temperatures specified on the boundaries and made from a material with constant thermal conductivity $\kappa = \kappa_e = 1.0$.

$$T = T_0 = 100^\circ \text{C} \quad \text{at} \quad r_0 = 20 \text{ mm} \quad (24)$$

$$T = T_n = 0^\circ \text{C} \quad \text{at} \quad r_n = 50 \text{ mm}$$

Consider two finite element models, each with four linear elements and five nodes: The first model (designated "Uniform Mesh") has equally spaced nodes. The second model (designated "Improved Mesh") has its nodal spacing governed by equation (23). The objective is the determination of the internal temperature distribution.

The solutions of the finite element governing equations lead to the results listed in table I. The error at the nodes is defined as the difference between the theoretical results and the finite element results. In order to compare the error distribution among the elements, rms errors were calculated using 50 uniformly spaced points through the thickness of each element. Each element's linear shape function was used to compute the temperature at any point along the element from nodal temperature values. Error values were calculated at each point from the difference between theoretical and finite element temperature values. The rms errors are shown beside the nodal errors in table I. The standard deviation of the within-element rms errors is shown at the bottom of table I with the overall rms error for the model which was calculated using 200 uniformly-spaced points through the model. (These are not quite the same points used for computing the within-element rms errors.)

The theoretical solution for this example (from eq. 3) is shown in figure 3. The nodal values for both models are superimposed on the theoretical solution. At this scale, differences between nodal and FEM values are much too small to be seen in the figure. The errors for the two models are plotted on an expanded scale in figure 4. Since linear elements were used, the errors tend to increase towards the center of each element (away from the nodes).

The uniform mesh has relatively small errors at the nodes and greater errors near the inner surface of the annular cylinder where the temperature gradient is greatest. The improved mesh is found to have zero errors at the nodes and a uniform distribution of errors between elements. The overall rms error of the improved mesh is 23 percent less than that of the uniform mesh.

Example 2

Next, consider the same cylinder but let the temperature gradient be specified on the inner boundary. Specifically, let the boundary conditions be:

$$\begin{aligned} \frac{dT}{dr} &= -5.4567833 \text{ } ^\circ\text{C/mm} \quad \text{at} \quad r_o = 20 \text{ mm} \\ T &= T_n = 0 \text{ } ^\circ\text{C} \quad \text{at} \quad r_n = 50 \text{ mm} \end{aligned} \tag{25}$$

(The boundary conditions were chosen to yield a nearly identical theoretical temperature distribution as in the first example.)

Table II shows the comparison between the finite element solution and the theoretical values of the temperature. Once again the improved mesh shows a more uniform distribution of rms errors between elements and the overall rms error is lower (by 15 percent) than that of the uniform mesh. The errors for the two models are plotted in figure 5.

Example 3

Since the temperature gradient is specified at the inner boundary, it is also of interest to know how the values of the temperature gradients obtained using the two finite element meshes compare with each other and with the theoretical values. Table III provides such a comparison. The errors as plotted in figure 6 are much greater (relatively) than in the previous examples, especially at the nodes. Note that the largest error is at the inside radius where the derivative boundary condition is applied. Generally the Neumann boundary conditions are not strongly satisfied in the finite element method (ref. 5). Therefore, in general, the FE method is not well suited for gradient-type problems. The improved mesh shows a slightly more uniform distribution of rms errors between elements and the overall rms error is also slightly lower (by 7 percent).

DISCUSSION

The numerical examples show that the "improved" mesh defined by equation (23) produces results which are in better agreement with the theoretical values than those obtained using the uniform mesh. Therefore, the mesh defined by equation (12) is an improvement over the uniform mesh for both the temperature-fixed boundary conditions and for the mixed boundary conditions.

In Example 1, the values of the stiffness matrix traces of the uniform and improved meshes are 74.6667 and 70.1520, respectively. This indicates that for this example the trace is only moderately sensitive to the nodal locations. More dramatic differences in the results will occur in problems where the trace is more sensitive to changes in the nodal positions. This will occur in problems with less regular geometry.

In the conventional finite element method, if a model with n degrees of freedom is refined by adding nodes, a new stiffness matrix needs to be constructed. In general the new stiffness matrix is different in form from the original stiffness matrix. However, in the "hierarchical" approach, new elements are constructed so that the original $n \times n$ stiffness matrix forms a submatrix of the new stiffness matrix. In this case, therefore, it is necessary to compute only the stiffness entries corresponding to the new degrees of freedom. This reduces the computing time in assembly. (See Zienkiewicz (ref. 6) for a more detailed discussion on hierarchical finite elements.) Since error estimation in the hierarchical approach has been established, we can take advantage of the error analyses in studying conventional finite elements.

Consider a finite element model with n degrees of freedom (d.o.f.). If the mesh is to be refined by introducing additional nodes (h-type refinement), then it is necessary to know the expected improvement in error before a refinement step is undertaken. O.C. Zienkiewicz, et al. (ref. 7) and Peano, et al. (ref. 8) have shown that if the $n+1$ th d.o.f. is to be introduced hierarchically, then the error in the energy norm is:

$$||e_{n,1}||^2 = \frac{(f_{n+1} - K_{n+1,n} u_n)^2}{K_{n+1,n+1}} \quad (26)$$

where f_{n+1} is the force corresponding to the $n+1$ th d.o.f., $K_{n+1,n+1}$ is the stiffness of the $n+1$ th d.o.f. $K_{n+1,n}$ is the off-diagonal stiffness relating the $n+1$ th d.o.f. to the original n d.o.f. system, and u_n is the array of nodal displacements of the n d.o.f. system. The subscripts $n,1$ of the error e refer to the n original d.o.f. and the new d.o.f.

Zienkiewicz (ref. 6) has used the above error relation to define an error indicator in the form:

$$\eta_{n,1}^2 = \frac{\left(\int_{\Omega} \zeta N_{n+1} d\Omega \right)^2}{K_{n+1,n+1}} \quad (27)$$

where ζ is the finite element residual.

In an adaptive refinement strategy, these indicators are normally calculated for all the d.o.f. corresponding to the next refinement. The indicators serve the purpose of identifying the region where refinement is necessary.

Next, the error corresponding to the previous iteration wherein the n th d.o.f. was added is:

$$||e_{n-1,1}||^2 = \frac{(f_n - K_{n,n-1} u_{n-1})^2}{K_{n,n}} \quad (28)$$

The corresponding error indicator is:

$$\eta_{n-1,1}^2 = \frac{\left(\int_{\Omega} \zeta N_n d\Omega \right)^2}{K_{n,n}} \quad (29)$$

These derivations are for hierarchical finite elements. However, the error with conventional finite elements will be of similar form (ref. 6).

The most general method of generating a good grid is to have an equal distribution of some specified weight function. (See Eiseman (ref. 9) for a complete discussion on adaptive grid generation.) Often, the error in the finite element solution is used as the weight function (ref. 10). Therefore the objective is to distribute the error equally among all elements. However, the value of the residual ζ can be obtained only after the equilibrium equations are solved. Nevertheless, one way of obtaining an equi-distribution of error a priori is by having a uniform element stiffness. As a consequence ζ will be nearly uniform among the elements. The trace minimization procedure developed herein produces such a result. Note that each of the ratios in the optimality condition, equation (19), is a constant γ .

$$\frac{r_1}{r_0} = \frac{r_2}{r_1} = \dots = \frac{r_{e+1}}{r_e} = \dots = \frac{r_n}{r_{n-1}} = \gamma \quad (30)$$

Substituting into equation (9), the element stiffness coefficient is:

$$S_e = \pi \kappa_e \frac{\gamma + 1}{\gamma - 1} \quad (31)$$

which is a constant. Therefore the trace minimization procedure produces a uniform element stiffness.

Finally, observe the graphs of the errors in figures 4 to 6. As mentioned earlier, the errors are more equally distributed with the improved mesh while there is a skewed distribution with the uniform mesh. In tables I to III, the standard deviation of the within-element rms errors is lower for the improved mesh than it is for the uniform mesh. Therefore, in all cases, the improved mesh distributes the error more uniformly than the uniform mesh. The rms errors are exactly equal (within precision of calculation) in the case where temperatures are specified at the boundaries. Therefore the mesh obtained in this case is optimal. Similar results, however are not obtained in the case where both temperature and gradient of temperature are specified because of the inability of FEM to strongly satisfy the Neumann boundary conditions. Nevertheless, it demonstrates the usefulness of the trace minimization procedure in a priori grid refinement.

CONCLUSIONS

The new method of finite-element grid improvement based on the minimization of the trace of the stiffness method has been extended to the problem of heat transfer in a solid body. In elasticity problems, this procedure is equivalent to minimizing the potential energy of the model by dividing the strain energy equally among the elements. The following conclusions were made:

1. Nodal positioning obtained by minimizing the trace of the stiffness matrix leads to an improved mesh over that obtained by uniform positioning of the nodes.
2. Since trace minimization is an a priori method, the mesh may be refined without solving the finite element problem. This makes the minimization procedure computationally inexpensive to perform. The mesh resulting from trace minimization may be used as a starting mesh for other mesh refinement procedures such as element division or element enhancement (h-methods or p-methods).
3. The method does not give satisfactory results for analyses in which both temperature and gradient of temperature are specified boundary conditions due to the inability of the FEM to strongly satisfy Neumann boundary conditions.

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TABLE I. - COMPARISON BETWEEN UNIFORM MESH, IMPROVED MESH, AND THEORETICAL VALUES FOR TEMPERATURE SPECIFIED BOUNDARY CONDITION PROBLEM

Radius, mm	Temperature value, °C			Uniform mesh error		Improved mesh error	
	Uniform mesh	Improved mesh	Theory	At node	In element rms	At node	In element rms
20.	100.	100.	100.	0.	-----	0.0000	-----
25.1487	-----	75.0000	75.	-----	-----	0.0000	0.5171
27.5	65.3550	-----	65.2453	-0.1096	1.0471	-----	-----
31.6228	-----	50.0000	50.	-----	-----	0.0000	0.5171
35.	39.0247	-----	38.9260	-0.0988	0.6688	-----	-----
39.7635	-----	25.0000	25.	-----	-----	0.0000	0.5171
42.5	17.7907	-----	17.7366	-0.0541	0.4422	-----	-----
50.	0.	0.	0.	0.	0.2857	0.0000	0.5171
Standard deviation of within-element rms errors					0.3309	0.00001	
Overall rms error					0.6801	0.5210	

TABLE II. - COMPARISON BETWEEN UNIFORM MESH, IMPROVED MESH, AND THEORETICAL VALUES FOR TEMPERATURE/TEMPERATURE-GRADIENT SPECIFIED BOUNDARY CONDITIONS

Radius, mm	Temperature value, °C			Uniform mesh error		Improved mesh error	
	Uniform mesh	Improved mesh	Theory	At node	In element rms	At node	In element rms
20.	99.4772	99.5700	100.	0.5228	-----	0.4300	-----
25.1487	-----	74.6775	75.	-----	-----	0.3225	0.2403
27.5	65.0133	-----	65.2453	0.2321	0.6786	-----	-----
31.6228	-----	49.7850	50.	-----	-----	0.2150	0.2980
35.	38.8207	-----	38.9260	0.1053	0.4273	-----	-----
39.7635	-----	24.8925	25.	-----	-----	0.1075	0.3782
42.5	17.6977	-----	17.7366	0.0389	0.3082	-----	-----
50.	0.	0.	0.	0.	0.2430	0.0000	0.4694
Standard deviation of within-element rms errors					0.1920	0.0996	
Overall rms error					0.4494	0.3803	

TABLE III. - COMPARISON BETWEEN UNIFORM MESH, IMPROVED MESH, AND THEORETICAL TEMPERATURE GRADIENT VALUES FOR TEMPERATURE/TEMPERATURE GRADIENT SPECIFIED BOUNDARY CONDITIONS

Radius, mm	Temperature gradient, °C/mm			Uniform mesh error		Improved mesh error	
	Uniform mesh	Improved mesh	Theory	At node	In element rms	At node	In element rms
20.	-4.5952	-4.8347	-5.4568	0.8616	-----	0.6220	-----
25.1487	-----	-3.84490	-4.3396	-----	-----	-0.4947	0.5183
27.5	-3.4923	-----	-3.9686	0.4762	0.6029	-----	-----
31.6228	-----	-3.0578	-3.4512	-----	-----	-0.3934	0.4122
35.	-2.8164	-----	-3.1182	0.3018	0.3594	-----	-----
39.7635	-----	-2.4317	-2.7446	-----	-----	-0.3129	0.3278
42.5	-2.3597	-----	-2.5679	0.2082	0.2392	-----	-----
50.	-2.3597	-2.4317	-2.1827	-0.1770	0.1134	-0.2490	0.1653
Standard deviation of within-element rms errors					0.2086	0.1490	
Overall rms error					0.3755	0.3478	

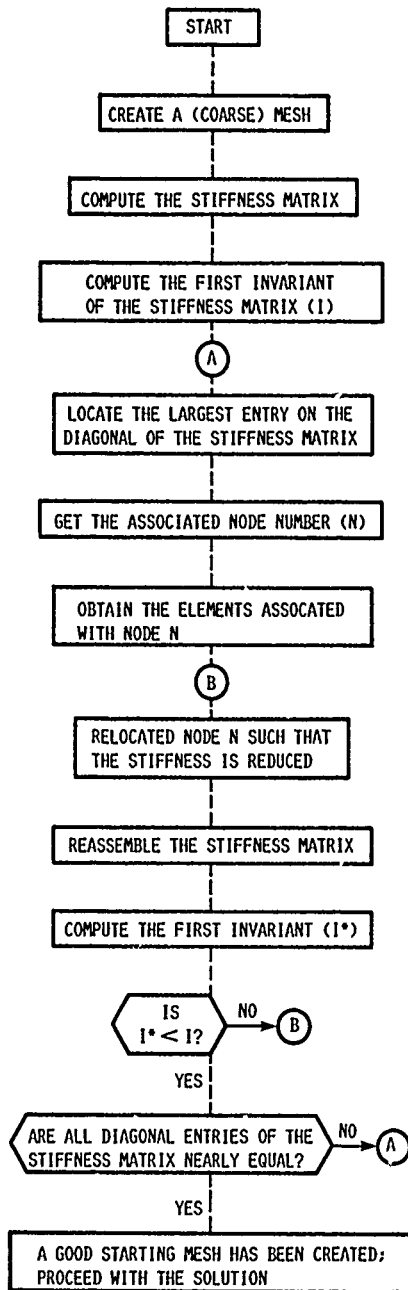


FIGURE 1. - FLOWCHART FOR MESH IMPROVEMENT PROCEDURE.

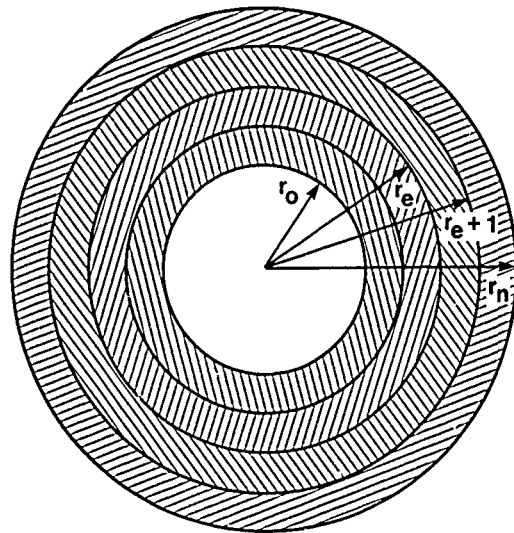


FIGURE 2. - FINITE ELEMENT MODEL.

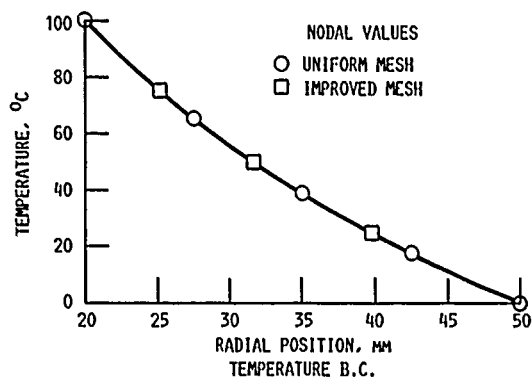


FIGURE 3. - SAMPLE PROBLEM: THEORETICAL AND NODAL TEMPERATURE VALUES.

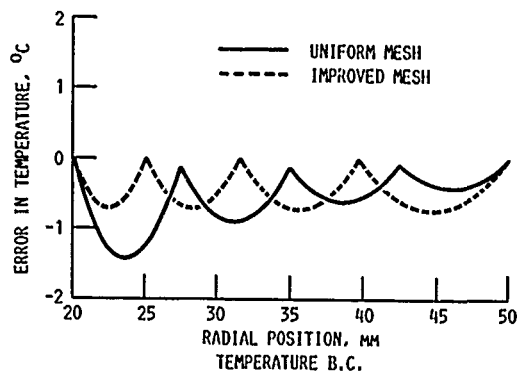


FIGURE 4. - EXAMPLE 1 (TEMPERATURE/TEMPORARY BOUNDARY CONDITIONS): ERRORS IN FEM RESULTS FOR CYLINDER TEMPERATURE.

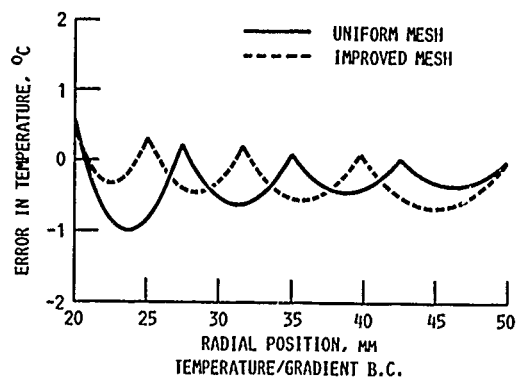


FIGURE 5. - EXAMPLE 2 (GRADIENT/TEMPERATURE BOUNDARY CONDITIONS): ERRORS IN FEM RESULTS FOR CYLINDER TEMPERATURE.

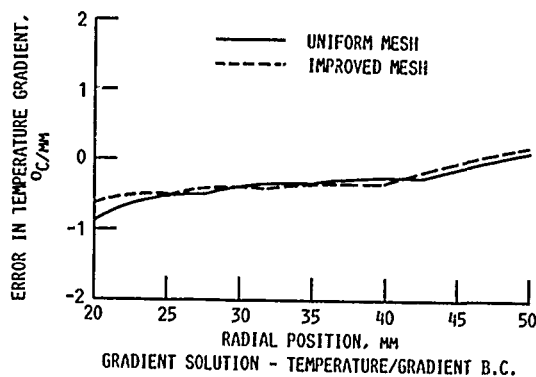


FIGURE 6. - EXAMPLE 2 (GRADIENT/TEMPERATURE BOUNDARY CONDITIONS): ERRORS IN FEM RESULTS FOR CYLINDER TEMPERATURE GRADIENT.

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16. Abstract This paper presents an extension of a structural finite element mesh improvement technique to heat conduction analysis. The mesh improvement concept was originally presented by Prager in studying tapered, axially loaded bars. It was further developed by Kittur, et al., who showed that an improved mesh can be obtained by minimizing the trace of the stiffness matrix. In this paper these procedures are extended and applied to the analysis of heat conduction in an infinitely long hollow circular cylinder.					
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